

NLO corrections to Z production in association with several jets

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in collaboration with

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- I: The numerical method**
- II: General improvements**
- III: First results for $pp \rightarrow Z + 5$ jets**

Part I

The numerical method

Z plus jet production at the LHC

Experimental status:

- The LHC experiments have measured Z production in association with up to 7 jets.

Theoretical status:

- NLO corrections to $Z + 0$ jets, $Z + 1$ jet, $Z + 2$ jets known for a long time.
- NLO corrections to $Z + 3$ jets and $Z + 4$ jets calculated by Blackhat collaboration.

Challenge:

- Can one calculate the NLO corrections to $Z + 5$ jets, $Z + 6$ jets and $Z + 7$ jets ?

Scaling behaviour with the number of jets

As the number of jets increases, the scaling behaviour with the number of jets is the relevant quantity.

- **Bad**: Factorial or exponential growth.
- **Better**: Polynomial growth.

Using recurrence relations, we can achieve n^3 -behaviour at LO.

What about loops ?

- Unitarity methods: n^9
- Numerical methods: n^3

Recurrence relations

Off-shell currents provide an efficient way to calculate amplitudes:

off-shell

$$= \sum_{j=1}^{n-1} \text{diagram}_1 + \sum_{j=1}^{n-2} \sum_{k=j+1}^{n-1} \text{diagram}_2$$

No Feynman diagrams are calculated in this approach !

The one-loop recurrence relations

$$\begin{aligned}
 & \text{Diagram 1} = \sum_{i=m}^{n-1} \text{Diagram 2} + \sum_{i=m}^{n-1} \text{Diagram 3} + \text{Diagram 4} \\
 & + \sum_{i=m}^{n-2} \sum_{j=i+1}^{n-1} \text{Diagram 5} + \sum_{i=m}^{n-2} \sum_{j=i+1}^{n-1} \text{Diagram 6} + \sum_{i=m}^{n-2} \sum_{j=i+1}^{n-1} \text{Diagram 7} \\
 & + \sum_{i=m}^{n-1} \text{Diagram 8} + \sum_{i=m}^{n-1} \text{Diagram 9} + \text{Diagram 10} + \text{Diagram 11}
 \end{aligned}$$

Numerical NLO QCD calculations

Proceed through the following steps:

1. **Local subtraction terms** for soft, collinear and ultraviolet singular part of the integrand of one-loop amplitudes
2. **Contour deformation** for the 4-dimensional loop integral.
3. **Numerical Monte Carlo integration** over phase space and loop momentum.

Not a new idea: Nagy and Soper proposed in '03 this method, working graph by graph.
(see also: Soper; Krämer, Soper; Catani et al.; Kilian, Kleinschmidt)

What is new: The IR-subtraction terms can be **formulated at the level of amplitudes**, no need to work graph by graph.

The IR-subtraction terms are **universal and amazingly simple**.

Subtraction method for loop integrals

Use subtraction also for the virtual part:

$$\int_{n+1} d\sigma^R + \int_n d\sigma^V = \underbrace{\int_{n+1} (d\sigma^R - d\sigma^A)}_{\text{convergent}} + \underbrace{\int_n (\mathbf{I} + \mathbf{L}) \otimes d\sigma^B}_n_{\text{finite}} + \underbrace{\int_n (d\sigma^V - d\sigma^{A'})}_{\text{convergent}}$$

- In the last term $d\sigma^V - d\sigma^{A'}$ the **Monte Carlo integration** is over a phase space integral of n final state particles plus a 4-dimensional loop integral.
- All **explicit poles cancel** in the combination $\mathbf{I} + \mathbf{L}$.
- Divergences of one-loop amplitudes related to **IR-divergences (soft and collinear)** and to **UV-divergences**.

The infrared subtraction terms for the virtual corrections

Local unintegrated form:

$$G_{\text{soft+coll}}^{(1)} = -4\pi\alpha_s i \sum_{i \in I_g} \left(\frac{4p_i p_{i+1}}{k_{i-1}^2 k_i^2 k_{i+1}^2} - 2 \frac{S_i g_{i-1,i}^{UV}}{k_{i-1}^2 k_i^2} - 2 \frac{S_{i+1} g_{i,i+1}^{UV}}{k_i^2 k_{i+1}^2} \right) A_i^{(0)}.$$

with $S_q = 1$, $S_g = 1/2$. The function $g_{i,j}^{UV}$ provides damping in the UV-region:

$$\lim_{k \rightarrow \infty} g_{i,j}^{UV} = O(k^{-2}), \quad \lim_{k_i \parallel k_j} g_{i,j}^{UV} = 1.$$

Integrated form:

$$S_\varepsilon^{-1} \mu^{2\varepsilon} \int \frac{d^D k}{(2\pi)^D} G_{\text{soft+coll}}^{(1)} = \frac{\alpha_s}{4\pi \Gamma(1-\varepsilon)} \sum_{i \in I_g} \left[\frac{2}{\varepsilon^2} \left(\frac{-2p_i \cdot p_{i+1}}{\mu^2} \right)^{-\varepsilon} + \frac{2}{\varepsilon} (S_i + S_{i+1}) \left(\frac{\mu_{\text{UV}}^2}{\mu^2} \right)^{-\varepsilon} \right] A_i^{(0)} + O(\varepsilon),$$

UV-subtraction terms

In a fixed direction in loop momentum space the **amplitude has up to quadratic UV-divergences**.

Only the **integration over the angles** reduces this to a logarithmic divergence.

For a local subtraction term we **have to match the quadratic, linear and logarithmic divergence**.

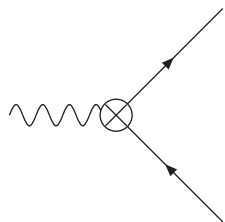
The subtraction terms have the **form of counter-terms** for propagators and vertices.

The complete UV-subtraction term **can be calculated recursively**.

UV-subtraction terms

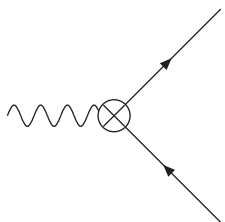
Example: **The quark-gluon vertex.**

Local unintegrated form:



$$= ig^3 S_\varepsilon^{-1} \mu^{4-D} \int \frac{d^D k}{(2\pi)^D i} \frac{2(1-\varepsilon) \bar{k} \not{\gamma}^\mu \bar{k} + 4\mu_{UV}^2 \gamma^\mu}{(\bar{k}^2 - \mu_{UV}^2)^3}$$

Integrated form:



$$= i \frac{g^3}{(4\pi)^3} \gamma^\mu (-1) \left(\frac{1}{\varepsilon} - \ln \frac{\mu_{UV}^2}{\mu^2} \right) + O(\varepsilon)$$

We can ensure that the integrated expression is proportional to the Born.

Contour deformation

With the subtraction terms for UV- and IR-singularities one removes

- UV divergences
- Pinch singularities due to **soft** or **collinear** partons

Still remains:

- **Singularities** in the integrand, **where a deformation** into the complex plane **of the contour is possible**.
- **Pinch singularities for exceptional configurations of the external momenta** (thresholds, anomalous thresholds ...)

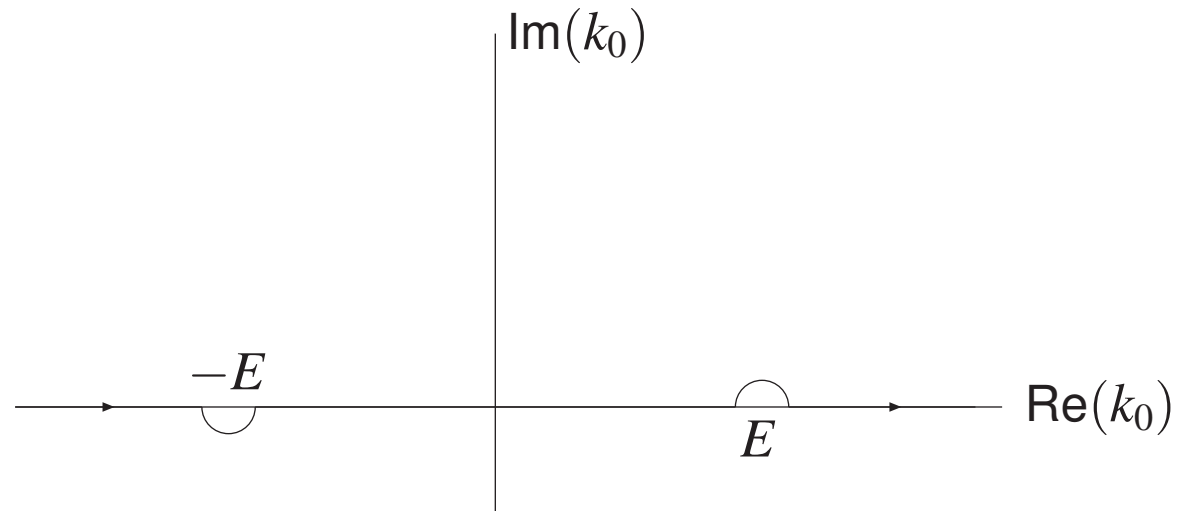
Contour deformation

$$I = \int d^4k \underbrace{[f(k) - g(k)]}_{h(k)}$$

$h(k)$ meromorphic function of four complex variables k_0, k_1, k_2, k_3 .

Integration over a surface of (real) dimension 4 in \mathbb{C}^4 .

I independent of the choice of the surface, as long as no poles are crossed.

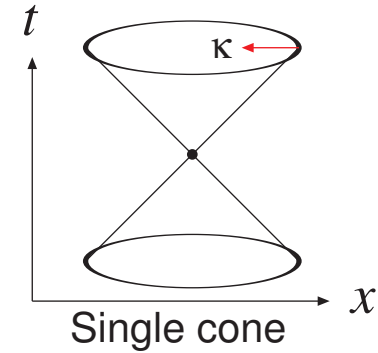


What is the best choice for the surface, in order to minimize Monte Carlo integration errors ?

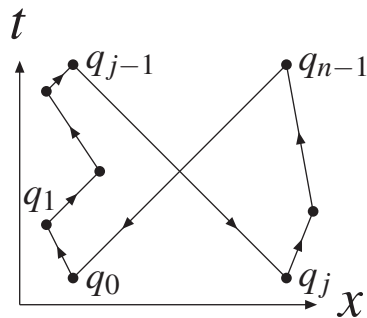
Direct contour deformation

Deformation of the loop momentum:

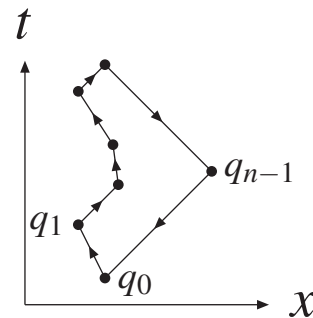
$$k_{\mathbb{C}} = k_{\mathbb{R}} + i\kappa$$



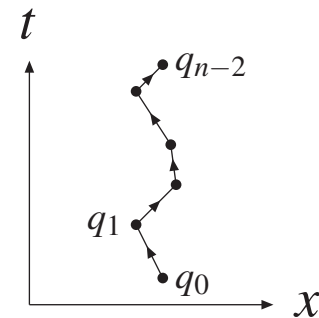
For n cones **draw only the origins** of the cones:



generic with 2 initial partons



initial partons adjacent



no initial partons

Efficiency

With the **local subtraction terms** and the **contour deformation** we obtain an integral, where the loop integration can – in principle – be performed with Monte Carlo methods.

However, the **integrand is oscillating**:

$$I = \int_0^1 dx [c + A \sin(2\pi x)], \quad A \gg c$$

This leads to **large Monte Carlo integration errors**.

Solution: **Antithetic variates**: Evaluate the integrand at x and $(1 - x)$.

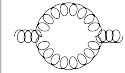
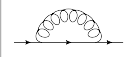
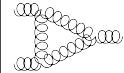
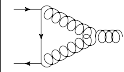
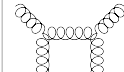
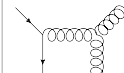
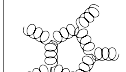
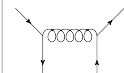
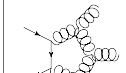
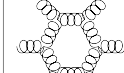
UV improvement

Ultraviolet behaviour of some example diagrams:

To the right: number of external particles

In the vertical:

leading power of the large $|k|$ -behaviour

	2	3	4	5	6
-2					
-3					
-4					
-5					
-6					

UV-finiteness requires fall off like $|k|^{-5}$.

$|k|^{-5}$ contribution is odd under $k \rightarrow -k$ and integrates to zero.

However, $|k|^{-5}$ term gives a large contribution to the Monte Carlo error.

UV improvement

- Split the integration **holomorphic** into two channels:

$$1 = \left[\prod_{j=1}^n \frac{k_j^2 - m_j^2}{\bar{k}^2 - \mu_{\text{UV}}^2} \right] + \left[1 - \prod_{j=1}^n \frac{k_j^2 - m_j^2}{\bar{k}^2 - \mu_{\text{UV}}^2} \right]$$

First channel: simple pole structure, can be evaluated with a simple contour.

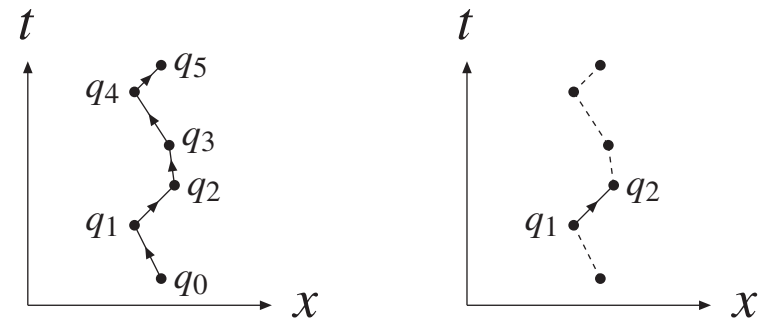
Second channel: Integrand falls off with two additional powers of $|k|$ in the ultraviolet.

- **Improvement of the counterterms** for the propagators and three-valent vertices from $|k|^{-5}$ to $|k|^{-7}$.
- Use **antithetic Monte Carlo integration technique**: Evaluate k and $(-k)$ together.

Infrared channels

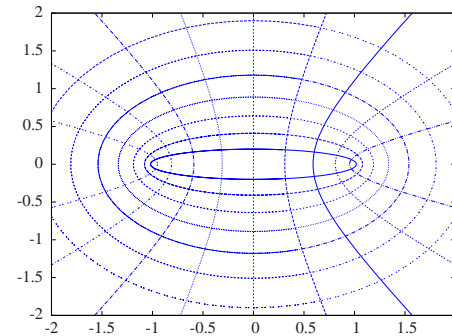
Non-holomorphic splitting:

$$I_{\text{int}} = \sum_i \int \frac{d^4 k}{(2\pi)^4} w_i(k) f(k),$$



Weights:

$$w_i(k) = \frac{\left(\frac{1}{|k_i^2| |k_{i+1}^2|} \right)^\alpha}{\sum_j \left(\frac{1}{|k_j^2| |k_{j+1}^2|} \right)^\alpha},$$



Coordinate system, where a line segment $[q_i, q_{i+1}]$ is singled out:

Generalised elliptical coordinates

Use technique of **antithetic variates** in these coordinates.

Part II

General improvements

Random polarisations

Matrix element with n external particles: **Instead of summing over all 2^n spin states**, introduce

$$\epsilon_\mu(\phi) = e^{i\phi}\epsilon_\mu^+ + e^{-i\phi}\epsilon_\mu^-.$$

and **replace the summation** over the spin states **by an integration** over the angle ϕ : P.

Draggiotis, R. Kleiss, C. Papadopoulos, '98

$$\sum_{\lambda=\pm} \epsilon_\mu^{\lambda*} \epsilon_\nu^\lambda = \frac{1}{2\pi} \int_0^{2\pi} d\phi \epsilon_\mu(\phi)^* \epsilon_\nu(\phi)$$

Works for **Born** and **virtual part** straightforward. For the **real part** the subtraction terms are usually spin-summed and thus **non-local** in ϕ .

Extension of the dipole formalism to random polarisations:

D. Götz, Ch. Schwan and S.W., '12

Colour decomposition at one-loop

One-loop amplitudes (and Born amplitudes with multiple quark pairs):

Partial amplitudes can be **decomposed** further **into primitive amplitudes** (gauge-invariant, cyclic ordered, fixed routing of fermions).

Z. Bern, L. Dixon, D. Kosower, '95

For amplitudes with more than one quark-antiquark pair this **decomposition is non-trivial**.

- Use Feynman diagrams and solve a (large) system of linear equations.

Ellis et al., '11; Ita, Ozeren, '11 ; Badger et al., '12

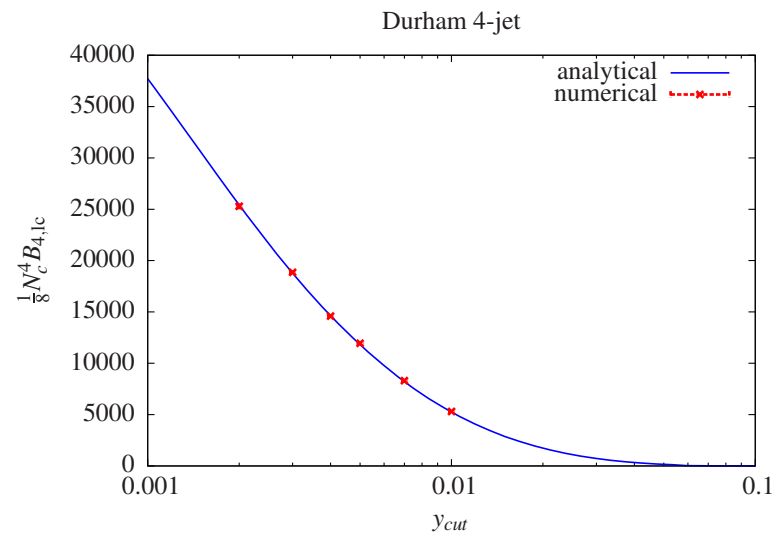
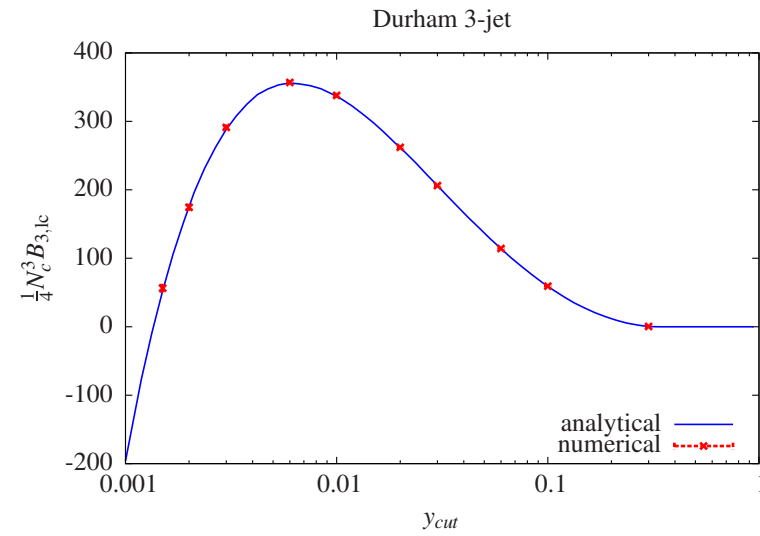
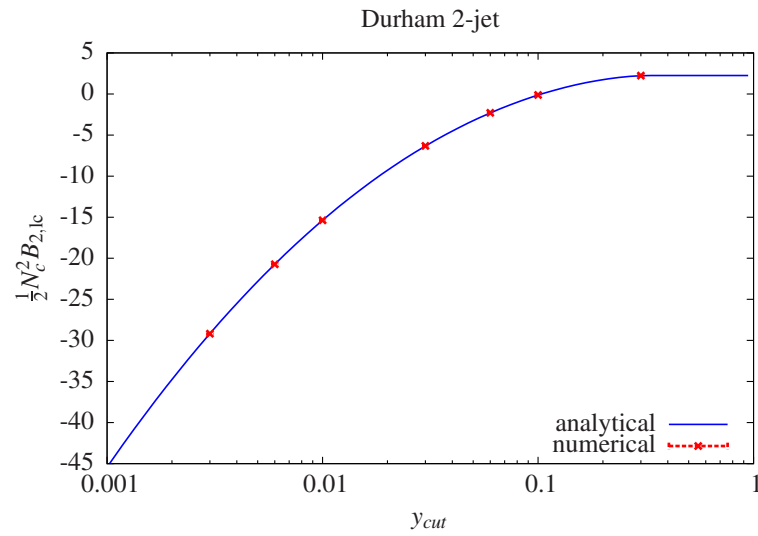
- More elegant: **Obtain colour decomposition directly** through shuffle relations.

Ch. Reuschle and S.W., '13

Part III

Numerical results

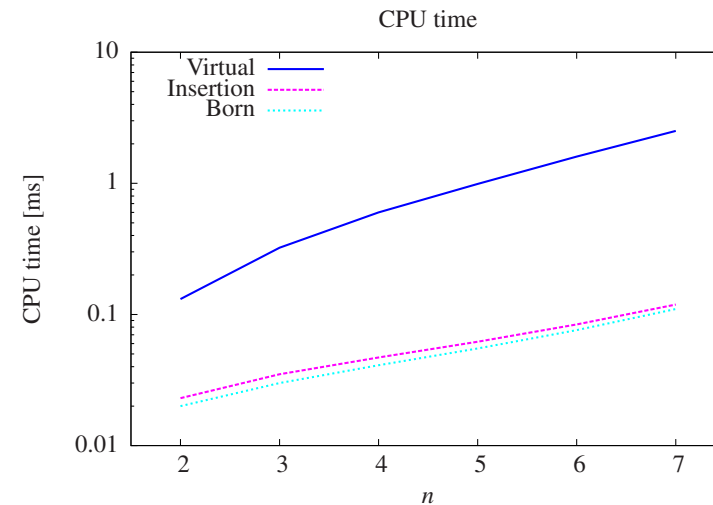
Jet rates in electron-positron annihilation



CPU scaling behaviour

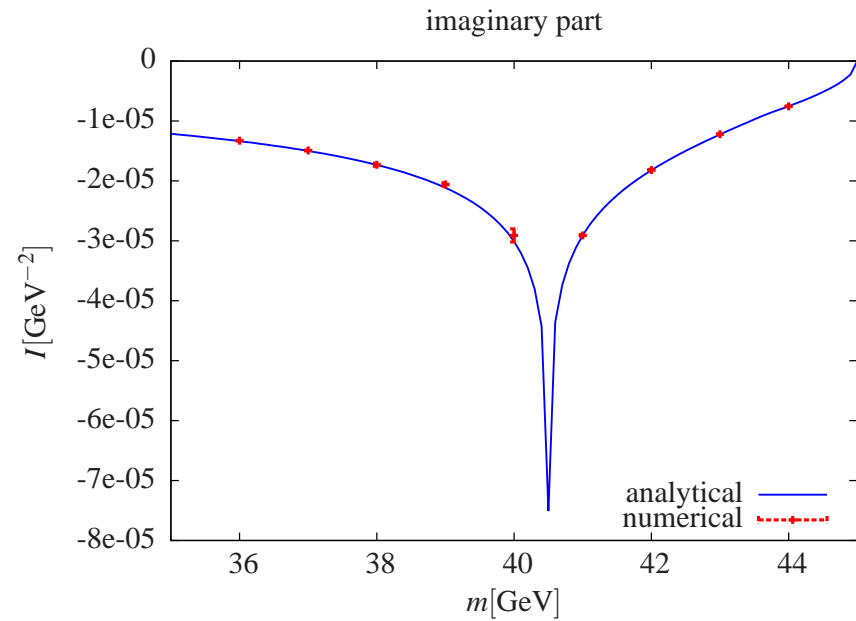
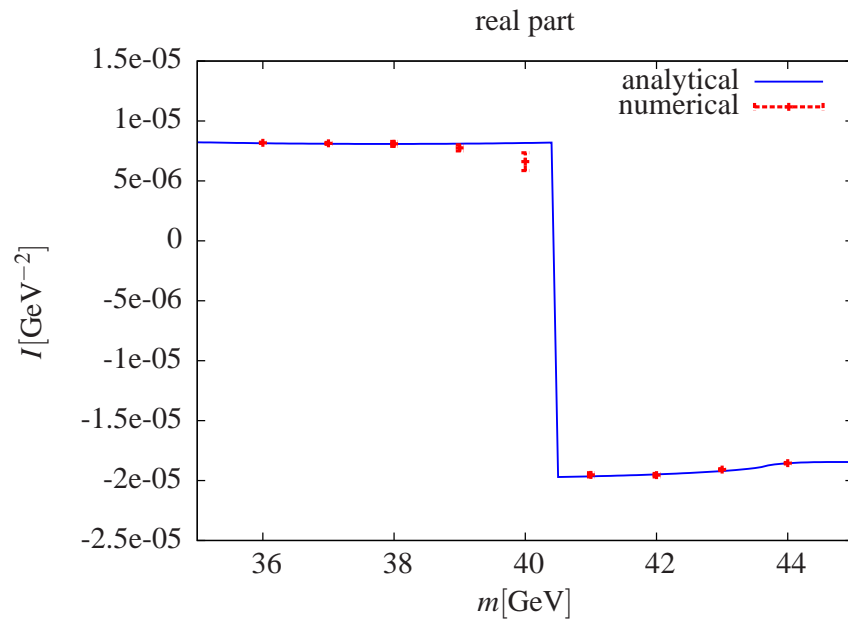
Scaling of the CPU time for one evaluation of the integrand (Born, virtual, insertion) with the number of external particles:

$$\text{CPU time} \sim n^4$$



- n^4 -behaviour from recurrence relations
- helicity summation replaced by smooth integration over random polarisations
- **Real part:** Extension of the dipole formalism to random polarisations

Extension to massive particles



Comparison of the results obtained by Monte Carlo integration with the analytical results in the vicinity of a threshold.

S. Becker and S.W., '12

Preliminary results on $pp \rightarrow Z + 5 \text{ jets}$

Process $pp \rightarrow Z + 5 \text{ jets} \rightarrow e^+e + 5 \text{ jets}$ at $\sqrt{s} = 7 \text{ TeV}$ with CTEQ6M/CTEQ6L1.
Jets defined by anti-kt-algorithm with $R = 0.5$.

Cuts:

$$p_l^\perp > 20 \text{ GeV}, \quad |\eta_l| < 2.5, \quad 66 \text{ GeV} < m_{l\bar{l}} < 116 \text{ GeV},$$
$$p_{\text{jet}}^\perp > 25 \text{ GeV}, \quad |\eta_{\text{jet}}| < 3.$$

Scale chosen on a per-event basis:

$$\mu_R = \mu_F = \frac{1}{2} H^{\perp'} = \frac{1}{2} \left(E_Z^\perp + \sum_j p_j^\perp \right).$$

Leading-colour approximation:

$$\sigma_{\text{LO,lc}} = 0.138 \pm 0.009 \text{ pb}, \quad \sigma_{\text{NLO,lc}} = 0.161 \pm 0.113 \text{ pb}.$$

Conclusions

- The numerical method for the computation of NLO corrections offers a good scaling behaviour.
- First results on $pp \rightarrow Z + 5$ jets.
- Public program available soon.